

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011104.D
 Acq On : 26 Jul 2022 11:36
 Operator : CG/JU
 Sample : PB146366BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK366

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/27/2022
 Supervised By :mohammad ahmed 07/27/2022

Quant Time: Jul 26 14:41:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 26 14:31:36 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	396550	20.000	ng/u1	0.00
20) Naphthalene-d8	10.457	136	1825134	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.328	164	1082101	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	1912403	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	1190239	20.000	ng/u1	# 0.00
88) Perylene-d12	23.551	264	1167299	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	75432	6.858	ng/uL	0.00
4) Pyridine-d5	3.570	84	1045432	36.770	ng/u1	0.00
7) Phenol-d5	6.840	99	1211580	37.220	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	845259	40.382	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	1014445m	38.568	ng/u1	0.00
15) 4-Methylphenol-d8	8.381	113	1022276	39.229	ng/u1	0.00
21) Nitrobenzene-d5	8.828	128	522215	39.262	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	539043	40.957	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	873287	36.283	ng/u1	0.00
31) 4-Chloroaniline-d4	10.604	131	1462921	36.869	ng/u1	0.00
46) Dimethylphthalate-d6	13.745	166	3123748	42.237	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	3608128	41.349	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	411827	27.514	ng/u1	0.00
60) Fluorene-d10	15.328	176	2446538	38.918	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	337598	35.675	ng/u1	0.00
73) Anthracene-d10	17.192	188	3374540	41.507	ng/u1	0.00
81) Pyrene-d10	19.445	212	3114324	49.820	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	2394232	42.705	ng/u1	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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