

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011123.D
 Acq On : 27 Jul 2022 11:20
 Operator : CG/JU
 Sample : PB146441BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK441

Quant Time: Jul 28 00:01:08 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	237425	20.000	ng/u1	0.00
20) Naphthalene-d8	10.451	136	993194	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.322	164	518948	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	970422	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	902528	20.000	ng/u1	# 0.00
88) Perylene-d12	23.545	264	965760	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	47004	7.137	ng/uL	0.00
4) Pyridine-d5	3.570	84	602977	35.422	ng/u1	0.00
7) Phenol-d5	6.840	99	684873	35.141	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	509197	40.631	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	568956	36.129	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	542607	34.777	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	284487	39.305	ng/u1	0.00
24) 2-Nitrophenol-d4	9.540	143	266613	37.226	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	442458	33.782	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	730497	33.831	ng/u1	0.00
46) Dimethylphthalate-d6	13.745	166	1373301	38.719	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	1704930	40.741	ng/u1	0.00
54) 4-Nitrophenol-d4	14.540	143	174552	24.317	ng/u1	0.00
60) Fluorene-d10	15.322	176	1135235	37.655	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	139633	29.079	ng/u1	0.00
73) Anthracene-d10	17.192	188	1671349	40.513	ng/u1	0.00
81) Pyrene-d10	19.445	212	1853999	39.113	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.398	264	1926657	41.537	ng/u1	0.00

Target Compounds	Qvalue

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

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