

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011119.D
 Acq On : 26 Jul 2022 20:49
 Operator : CG/JU
 Sample : PB146439BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK439

Quant Time: Jul 26 22:52:13 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	389914	20.000	ng/u1	0.00
20) Naphthalene-d8	10.452	136	1709820	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.322	164	943101	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	1620611	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.204	240	1260261	20.000	ng/u1	#-0.01
88) Perylene-d12	23.545	264	1383851	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	68051	6.292	ng/uL	0.00
4) Pyridine-d5	3.570	84	914342	32.707	ng/u1	0.00
7) Phenol-d5	6.840	99	1065806	33.299	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	752391	36.557	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	887729	34.325	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	864465	33.737	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	460273	36.939	ng/u1	0.00
24) 2-Nitrophenol-d4	9.540	143	457495	37.106	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	725144	32.160	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	1186450	31.918	ng/u1	0.00
46) Dimethylphthalate-d6	13.746	166	2275007	35.295	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	2813118	36.989	ng/u1	0.00
54) 4-Nitrophenol-d4	14.540	143	287151	22.012	ng/u1	0.00
60) Fluorene-d10	15.322	176	1837581	33.539	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	222245	27.714	ng/u1	0.00
73) Anthracene-d10	17.192	188	2463817	35.761	ng/u1	0.00
81) Pyrene-d10	19.445	212	2461462	37.188	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.398	264	2354782	35.429	ng/u1	0.00

Target Compounds	Qvalue

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

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