

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011144.D
 Acq On : 28 Jul 2022 00:02
 Operator : CG/JU
 Sample : PB146455BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK455

Quant Time: Jul 28 01:07:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 00:46:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	446999	20.000	ng/u1	0.00
20) Naphthalene-d8	10.445	136	1871012	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.322	164	939465	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.086	188	1629679	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.204	240	1508916	20.000	ng/u1	# 0.00
88) Perylene-d12	23.539	264	1696009	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	58843	4.746	ng/uL	0.00
4) Pyridine-d5	3.569	84	814177	25.405	ng/u1	0.00
7) Phenol-d5	6.834	99	1077570	29.367	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	777450	32.950	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	886403	29.897	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	845918	28.798	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	436780	32.034	ng/u1	0.00
24) 2-Nitrophenol-d4	9.540	143	428252	31.741	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	691556	28.028	ng/u1	0.00
31) 4-Chloroaniline-d4	10.598	131	1098110	26.996	ng/u1	0.00
46) Dimethylphthalate-d6	13.739	166	1970321	30.686	ng/u1	0.00
49) Acenaphthylene-d8	14.010	160	2507397	33.097	ng/u1	0.00
54) 4-Nitrophenol-d4	14.534	143	249090	19.168	ng/u1	0.00
60) Fluorene-d10	15.322	176	1596840	29.258	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	190871	23.669	ng/u1	0.00
73) Anthracene-d10	17.186	188	2199462	31.747	ng/u1	0.00
81) Pyrene-d10	19.439	212	2425337	30.604	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.392	264	2649348	32.524	ng/u1	0.00

Target Compounds Qvalue

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

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