

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
 Data File : BP014653.D
 Acq On : 11 Apr 2023 13:26
 Operator : CG/JU
 Sample : 02251-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMC0

Quant Time: Apr 11 23:01:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 11 01:21:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	125920	20.000	ng/ul	0.00
20) Naphthalene-d8	10.793	136	543732	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	329639	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	637345	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	531053	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	654603	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	16035	4.938	ng/uL	0.00
4) Pyridine-d5	3.799	84	518	0.062	ng/ul	0.00
7) Phenol-d5	7.152	99	85128	7.330	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	233447	31.331	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	206995	24.661	ng/ul	0.00
15) 4-Methylphenol-d8	8.699	113	151398	16.162	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	120368	27.422	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	139696	27.793	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	215181	23.507	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	46608	3.866	ng/ul	0.02
46) Dimethylphthalate-d6	14.040	166	806535	30.176	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	883948	29.806	ng/ul	0.00
54) 4-Nitrophenol-d4	14.851	143	20	0.005	ng/ul	0.00
60) Fluorene-d10	15.622	176	664362	29.807	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	107162	22.851	ng/ul	0.00
73) Anthracene-d10	17.492	188	1054889	34.200	ng/ul	0.00
81) Pyrene-d10	19.722	212	1133929	31.789	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1211504	34.835	ng/ul	0.00

Target Compounds Qvalue

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

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